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CORRECTION

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Correction: Sumanene derivatives functionalized at the internal carbon

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Correction for 'Sumanene derivatives functionalized at the internal carbon' by Niti Ngamsomprasert *et al.*, *Chem. Commun.*, 2017, **53**, 697–700.

The authors regret that there was an error on page 698 of the original article. The text originally read, "The space group of the crystal was $P3_2$ and the antipode (+)-**5** gave the $P3_1$ system." This should have read, "The space group of the crystal was $P3_1$ and the antipode (+)-**5** gave the $P3_2$ system."

The Royal Society of Chemistry apologises for these errors and any consequent inconvenience to authors and readers.

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